

## Supporting Information

### Density functional theory study on silver and bis-silver complexes with lighter tetrylene – Are silver and bis-silver carbenes candidates for SARS-CoV-2 inhibition? An insight from molecular docking simulation

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## Table of Contents

Contents	Page
<b>Table S1.</b> Cartesian coordinates of the optimized ligands NHC <sub>ph</sub> to NHGe <sub>ph</sub> ( <b>NHC – NHGe</b> ) at the BP86/def2-SVP level of theory.	S2
<b>Table S2.</b> Cartesian coordinates of the optimized silver- tetrylene complexes AgCl-NHC <sub>ph</sub> to AgCl-NHGe <sub>ph</sub> ( <b>Ag-NHC – Ag-NHGe</b> ) at the BP86/def2-SVP level of theory.	S5
<b>Table S3.</b> Cartesian coordinates of the optimized bis-silver-tetrylene complexes (NHC <sub>ph</sub> -AgCl) <sub>2</sub> to (NHGe <sub>ph</sub> -AgCl) <sub>2</sub> ( <b>Ag-NHC-bis – Ag-NHGe-bis</b> ) at the BP86/def2-SVP level of theory.	S8
<b>Figure S1.</b> Optimized geometries of silver-complexes <b>NHE-Ag</b> (E = C, Si, Ge) at the BP86/def2-SVP level with dispersive effect. Bond lengths are given in Å; angles in degrees.	S13
<b>Figure S2.</b> Optimized geometries of silver-complexes <b>NHE-Ag-bis</b> (E = C, Si, Ge) at the BP86/def2-SVP level with dispersive effect. Bond lengths are given in Å; angles in degrees	S14